AMENDMENTS TO THE CLAIMS:

This listing of claims will replace all prior versions and listings of claims in the application:

1. (Currently Amended) A compound of the formula:

$$R^{5}$$
 R^{6}
 R^{6

wherein

ring A represents a benzene an aromatic ring optionally having substituents;

B <u>represents a C₁₋₆ alkylene optionally having substituents;</u> [[,]] Y and Ya are the same or different and each represents a bond or a spacer having a main chain of 1 to 6 atoms;

R¹ and R² are the same or different and each represents a hydrogen atom, a hydrocarbon group optionally having substituents or a heterocyclic group optionally having substituents, or R¹ and R², together with the adjacent nitrogen atom, form a nitrogen-containing heterocyclic ring optionally having substituents, or R¹ is linked with

ring A together with the adjacent nitrogen atom and B to form a 5- to 7-membered nitrogen-containing heterocyclic ring;

R³ represents a hydrogen atom, a hydrocarbon group optionally having substituents or a heterocyclic group optionally having substituents;

R⁴ and R⁵ are the same or different and each represents a hydrogen atom or a hydrocarbon group optionally having substituents, or R⁴ and R⁵, together with the adjacent carbon atom, form a ring optionally having substituents;

R⁶ represents an indolyl group optionally having substituents; and

Z represents piperidinyl optionally having substituents or piperazinyl optionally having substituents; and

Za are the same or different and each represents a hydrogen atom, a halogen atom or a cyclic group optionally having substituents; or a salt thereof.

- 2. (Canceled)
- 3. (Original) The compound according to claim 1, wherein R^3 is a hydrogen atom or a C_{1-6} alkyl optionally having substituents.
- 4. (Original) The compound according to claim 1, wherein one of R^4 and R^5 is a hydrogen atom, and the other is a C_{1-6} alkyl optionally having substituents.
 - 5-6. (Canceled)
- 7. (Currently Amended) The compound according to claim 1 [[5]], wherein Z is piperidinyl or piperazinyl, each of which is substituted by a group of the formula: -Yd-Ara wherein Yd represents a bond or a spacer having a main chain of 1 to 6 atoms, and Ara represents a monocyclic group optionally having substituents.

- 8. (Original) The compound according to claim 1, wherein Ya is a bond, and Za is a hydrogen atom.
 - 9. (Original) The compound according to claim 1, wherein B is a C₁₋₆ alkylene.10. (Canceled)
- 11. (Original) The compound according to claim 1, wherein R^1 and R^2 are C_{1-6} alkyl.
 - 12. (Original) The compound according to claim 1, wherein Y is -CO-.
 - 13. (Original) The compound according to claim 1, which is

N-((1R.2S)-1-(((5-((dimethylamino)methyl)-2-

((methylamino)carbonyl)phenyl)amino)carbonyl)-2-(1H-indol-3-yl)propyl)-4-(2-methylphenyl)-1-piperidinecarboxamide;

N-((1R,2S)-1-(((2-((dimethylamino)carbonyl)-5-((dimethylamino)methyl)phenyl)amino)carbonyl)-2-(1H-indol-3-yl)propyl)-4-(4-fluorophenyl)-1-piperidinecarboxamide;

N-((1R,2S)-1-(((5-((dimethylamino)methyl)-2-methoxyphenyl)amino)carbonyl)-2-(1H-indol-3-yl)propyl)-4-(4-fluoro-2-methylphenyl)-3-oxo-1-piperazinecarboxamide;

N-((1R,2S)-1-(((5-((dimethylamino)methyl)-2-methoxyphenyl)amino)carbonyl)-2-(1H-indol-3-yl)propyl)-4-(2-methylphenyl)-1-piperazinecarboxamide;

N-((1R,2S)-1-(((5-((dimethylamino)methyl)-2-ethoxyphenyl)amino)carbonyl)-2-(1H-indol-3-yl)propyl)-4-(4-fluorophenyl)-1-piperazinecarboxamide; or

N-((1R,2S)-1-(((5-((dimethylamino)methyl)-2-ethoxyphenyl)amino)carbonyl)-2-(1H-indol-3-yl)propyl)-4-phenyl-1-piperidinecarboxamide.

- 14. (Currently Amended) A pharmaceutical preparation comprising the compound according to claim 1[[,]] or a salt thereof or a prodrug thereof.
- 15. (Original) The pharmaceutical preparation according to claim 14, which is a somatostatin receptor binding inhibitor.
- 16. (Original) The pharmaceutical preparation according to claim 15, which is a somatostatin subtype 2 receptor binding inhibitor.
- 17. (Original) The pharmaceutical preparation according to claim 14, which is a somatostatin receptor agonist.
- 18. (Original) The pharmaceutical preparation according to claim 17, which is a somatostatin subtype 2 receptor agonist.
- 19. (Original) The pharmaceutical preparation according to claim 14, which is a prophylactic or therapeutic agent for diabetes or diabetic complications.
- 20. (Original) The pharmaceutical preparation according to claim 14, which is a prophylactic or therapeutic agent for obesity.
 - 21. (Canceled)
- 22. (Currently Amended) A method for inhibiting somatostatin receptor binding in a mammal, which comprises administering to the mammal an effective amount of the compound according to claim 1[[,]] or a salt thereof or a prodrug thereof.
 - 23. (Canceled)
- 24. (Currently Amended) A method for preventing or treating diabetes or diabetic complications in a mammal, which comprises administering to the mammal an effective amount of the compound according to claim 1[[,]] or a salt thereof—or a prodrug-thereof.

25. (Canceled)

- 26. (Currently Amended) A method for preventing or treating obesity in a mammal, which comprises administering to the mammal an effective amount of the compound according to claim 1[[,]] or a salt thereof or a prodrug thereof.
- 27. (Currently Amended) A method for producing a compound of claim 1 or a salt thereof, which comprises reacting a compound of the formula:

$$R^{5}$$
 R^{6}
 N
 N
 Y - Z

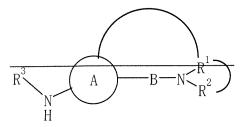
wherein

Y represents a bond or a spacer having a main chain of 1 to 6 atoms;

R⁴ and R⁵ are the same or different, and each represents a hydrogen atom or a hydrocarbon group optionally having substituents, or R⁴ and R⁵, together with the adjacent carbon atom, form a ring optionally having substituents;

R⁶ represents an indolyl group optionally having substituents;

Z represents <u>piperidinyl optionally having substituents or piperazinyl optionally having</u>
<u>substituents</u> a hydrogen atom, a halogen atom or a cyclic group optionally having
<u>substituents</u>; or a salt thereof, with a compound of the formula:



$$R^{3} \xrightarrow{M} H = N < R^{1}$$

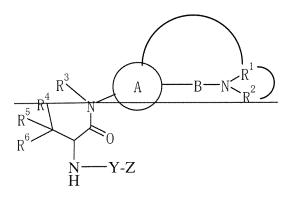
wherein

ring A represents <u>a benzene</u> an aromatic ring optionally having substituents;

B represents a C₁₋₆ alkylene optionally having substituents bond or a spacer having a main chain of 1 to 6 atoms;

R¹ and R² are the same or different, and each represents a hydrogen atom, a hydrocarbon group optionally having substituents or a heterocyclic group optionally having substituents, or R¹-and R²-, together with the adjacent nitrogen atom, form a nitrogen-containing heterocyclic ring optionally having substituents, or R¹-is linked with ring A together with the adjacent nitrogen atom and B to form a 5- to 7-membered nitrogen-containing heterocyclic ring;

R³ represents a hydrogen atom, a hydrocarbon group optionally having substituents or a heterocyclic group optionally having substituents; or a salt thereof to give a compound of the formula:



wherein

each symbol is as defined above; or a salt thereof, and optionally reacting the compound or a salt thereof with a compound of the formula: L⁴-Ya-Za wherein L⁴ represents a leaving group; Ya represents a bond or a spacer having a main chain of 1 to 6 atoms; Za represents a hydrogen atom, a halogen atom or a cyclic group optionally having substituents; or a salt thereof.

28. (Canceled)